

Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG120921\
 Data File : BG051467.D
 Acq On : 10 Dec 2021 23:05
 Operator : CG/JU
 Sample : PB141304BL
 Misc :
 ALS Vial : 16 Sample Multiplier: 1

Instrument :
 BNA_G
 ClientSampleId :
 SBLK304

Quant Time: Dec 11 04:26:08 2021
 Quant Method : Z:\svoasrv\HPCHEM1\BNA_G\Methods\SFAM-EPA-BG120821.M
 Quant Title : SVOA CALIBRATION
 QLast Update : Thu Dec 09 03:21:41 2021
 Response via : Initial Calibration

Compound	R.T.	QIon	Response	Conc	Units	Dev(Min)

Internal Standards						
1) 1,4-Dichlorobenzene-d4	8.188	152	26609	20.000	ng/u1	0.00
20) Naphthalene-d8	11.014	136	118720	20.000	ng/u1	0.00
38) Acenaphthene-d10	14.815	164	80424	20.000	ng/u1	0.00
64) Phenanthrene-d10	17.565	188	186463	20.000	ng/u1	0.00
79) Chrysene-d12	21.872	240	165922	20.000	ng/u1	0.00
88) Perylene-d12	25.262	264	147938	20.000	ng/u1	-0.01

System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.534	96	4867	6.007	ng/uL	0.00
4) Pyridine-d5	3.975	84	58771	25.258	ng/u1	0.01
7) Phenol-d5	7.365	99	82684	30.524	ng/u1	0.01
9) Bis-(2-Chloroethyl)eth...	7.500	67	54600	31.431	ng/u1	0.00
11) 2-Chlorophenol-d4	7.723	132	60412	31.346	ng/u1	0.00
15) 4-Methylphenol-d8	8.916	113	64780	30.441	ng/u1	0.00
21) Nitrobenzene-d5	9.369	128	32659	31.712	ng/u1	0.00
24) 2-Nitrophenol-d4	10.097	143	37362	32.060	ng/u1	0.00
28) 2,4-Dichlorophenol-d3	10.655	165	52946	27.927	ng/u1	0.00
31) 4-Chloroaniline-d4	11.167	131	83447	30.093	ng/u1	0.00
46) Dimethylphthalate-d6	14.216	166	214733	34.506	ng/u1	0.00
49) Acenaphthylene-d8	14.516	160	258546	32.804	ng/u1	0.00
54) 4-Nitrophenol-d4	15.097	143	23270	24.836	ng/u1	0.03
60) Fluorene-d10	15.808	176	181111	32.693	ng/u1	0.00
65) 4,6-Dinitro-2-methylph...	15.955	200	29934	27.015	ng/u1	0.00
73) Anthracene-d10	17.665	188	296165	33.946	ng/u1	0.00
81) Pyrene-d10	19.944	212	352556	35.351	ng/u1	0.00
92) Benzo(a)pyrene-d12	25.027	264	280699	36.787	ng/u1	0.00

Target Compounds Qvalue

(#) = qualifier out of range (m) = manual integration (+) = signals summed

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